LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090321\
Data File : VX024111.D
Acq On : 03 Sep 2021 18:37
Operator : JC/MD
Sample : M3646-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
1992

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Title : SW846 8260

Signal : TIC: VX024111.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.148	7	10	17	rVB2	7961	10982	1.38%	0.139%
2	1.362	42	45	51	rVB2	7223	8563	1.07%	0.108%
3	1.453	56	60	67	rBV	43437	50623	6.35%	0.641%
4	2.300	194	199	206	rVB	5314	9063	1.14%	0.115%
5	2.386	206	213	227	rVB	108136	178067	22.32%	2.255%
6	2.550	230	240	249	rBV	22341	41277	5.17%	0.523%
7	2.648	251	256	262	rVB	5005	8919	1.12%	0.113%
8	2.971	299	309	319	rVB4	8289	22248	2.79%	0.282%
9	4.580	561	573	584	rBV3	13418	42684	5.35%	0.541%
10	5.397	693	707	723	rBV	92602	270793	33.94%	3.430%
11	5.562	723	734	749	rVV	147600	413858	51.88%	5.241%
12	5.964	789	800	809	rBV	104330	273837	34.33%	3.468%
13	6.056	809	815	822	rVB2	18090	47705	5.98%	0.604%
14	6.208	832	840	847	rBV4	3147	8608	1.08%	0.109%
15	6.769	922	932	942	rBV	231707	548147	68.71%	6.942%
16	7.470	1041	1047	1055	rVB2	4531	9904	1.24%	0.125%
17	7.909	1112	1119	1125	rVB2	12151	24101	3.02%	0.305%
18	8.580	1221	1229	1234	rVV2	5754	11420	1.43%	0.145%
19	8.653	1234	1241	1248	rVV	516156	797765	100.00%	10.104%
20	8.720	1248	1252	1263	rVB	40381	66222	8.30%	0.839%
21	9.433	1362	1369	1378	rVB7	5055	10935	1.37%	0.138%
22	10.055	1462	1471	1475	rBV	502163	685836	85.97%	8.686%
23	10.092	1475	1477	1489	rVB	49759	66972	8.39%	0.848%
24	10.195	1489	1494	1500	rBV	20740	29625	3.71%	0.375%
25	10.305	1507	1512	1518	rVB	47071	61874	7.76%	0.784%
26	10.390	1518	1526	1529	rBV2	6697	11016	1.38%	0.140%
27	10.646	1563	1568	1573	rBV2	40993	59212	7.42%	0.750%
28	11.018	1625	1629	1634	rVB3	5865	9823	1.23%	0.124%
29	11.085	1634	1640	1651	rBV	455182	569891	71.44%	7.218%
30	11.177	1651	1655	1660	rVB4	4939	9608	1.20%	0.122%
31	11.384	1684	1689	1695	rBV	19149	33146	4.15%	0.420%
32	11.457	1697	1701	1707	rVB3	9720	13663	1.71%	0.173%
33	11.616	1722	1727	1740	rVB3	34346	62263	7.80%	0.789%
34	11.756	1744	1750	1755	rBV2	44877	61238	7.68%	0.776%
35	11.811	1755	1759	1762	rVV	28519	34668	4.35%	0.439%
36	11.841	1762	1764	1770	rVV2	11917	17925	2.25%	0.227%

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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : 1992

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Title : SW846 8260

37	11.976	1779	1786	1789	rVV5	12515	24757	3.10%	0.314%
38	12.024	1789	1794	1799	rVV	518988	641364	80.40%	8.123%
39	12.104	1802	1807	1819	rVB2	276425	397066	49.77%	5.029%
40	12.244	1826	1830	1834	rBV2	16679	22219	2.79%	0.281%
41	12.317	1839	1842	1846	rBV3	9577	12785	1.60%	0.162%
42	12.414	1853	1858	1862	rVV3	21185	30331	3.80%	0.384%
43	12.469	1862	1867	1872	rVB5	8494	15044	1.89%	0.191%
44	12.561	1878	1882	1886	rVB4	31000	44523	5.58%	0.564%
45	12.616	1886	1891	1894	rVB3	9624	13160	1.65%	0.167%
46	12.664	1894	1899	1901	rBV4	7346	11846	1.48%	0.150%
47	12.695	1901	1904	1907	rVV2	12863	16638	2.09%	0.211%
48	12.756	1911	1914	1919	rVV4	13047	24546	3.08%	0.311%
49	12.799	1919	1921	1927	rVB4	9747	10283	1.29%	0.130%
50	12.853	1927	1930	1934	rBV2	11501	14370	1.80%	0.182%
51	12.920	1935	1941	1944	rBV5	7615	14718	1.84%	0.186%
52	12.969	1944	1949	1953	rBV3	13390	22291	2.79%	0.282%
53	13.012	1953	1956	1960	rVB4	10855	14459	1.81%	0.183%
54	13.067	1960	1965	1972	rBV10	8553	19682	2.47%	0.249%
55	13.128	1972	1975	1978	rVV5	8786	13123	1.64%	0.166%
56	13.170	1979	1982	1987	rVB6	12674	17850	2.24%	0.226%
57	13.256	1992	1996	1999	rVV2	32367	41688	5.23%	0.528%
58	13.292	1999	2002	2007	rVB2	30100	40598	5.09%	0.514%
59	13.347	2007	2011	2013	rBV	27840	36361	4.56%	0.461%
60	13.372	2013	2015	2018	rVV	25536	23993	3.01%	0.304%
61	13.426	2018	2024	2029	rVB2	56985	86158	10.80%	1.091%
62	13.512	2033	2038	2047	rBV	197838	262971	32.96%	3.330%
63	13.591	2047	2051	2058	rVB8	12997	28766	3.61%	0.364%
64	13.664	2058	2063	2067	rBV4	10807	17888	2.24%	0.227%
65	13.731	2070	2074	2078	rBV2	30813	48402	6.07%	0.613%
66	13.780	2078	2082	2088	rVB	59700	89606	11.23%	1.135%
67	13.853	2090	2094	2097	rBV5	10351	17279	2.17%	0.219%
68	13.884	2097	2099	2103	rVV4	12432	14655	1.84%	0.186%
69	13.926	2104	2106	2111	rVB4	7056	9581	1.20%	0.121%
70	13.981	2111	2115	2117	rBV4	13142	19347	2.43%	0.245%
71	14.006	2117	2119	2122	rVB2	15027	15225	1.91%	0.193%
72	14.079	2127	2131	2133	rBV5	9390	12667	1.59%	0.160%
73	14.134	2137	2140	2145	rBV5	16770	19350	2.43%	0.245%
74	14.231	2150	2156	2159	rBV4	31995	55724	6.99%	0.706%
75	14.262	2159	2161	2167	rVB	19713	20436	2.56%	0.259%
76	14.323	2167	2171	2176	rBV3	24137	33987	4.26%	0.430%

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Misc   : 5.0mL/MSVOA_X/WATER
ALS Vial : 23   Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : 1992

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON   Filtering: 5
Sampling : 1   Min Area: 3 % of largest Peak
Start Thrs: 0.2   Max Peaks: 100
Stop Thrs : 0   Peak Location: TOP

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Method   : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Title   : SW846 8260

77	14.378	2176	2180	2183	rVB2	11877	13858	1.74%	0.176%
78	14.481	2192	2197	2201	rBV2	31256	47306	5.93%	0.599%
79	14.524	2201	2204	2210	rVB5	7727	11912	1.49%	0.151%
80	14.640	2220	2223	2226	rVB	49225	52736	6.61%	0.668%
81	14.676	2226	2229	2233	rVB3	19481	25104	3.15%	0.318%
82	14.725	2233	2237	2240	rBV5	10897	18327	2.30%	0.232%
83	14.780	2240	2246	2250	rBV2	56770	90903	11.39%	1.151%
84	14.853	2255	2258	2260	rVB4	9799	10443	1.31%	0.132%
85	14.908	2264	2267	2272	rVB7	8585	13329	1.67%	0.169%
86	15.048	2287	2290	2294	rVB6	6920	11421	1.43%	0.145%
87	15.188	2308	2313	2315	rBV	25944	40979	5.14%	0.519%
88	15.377	2340	2344	2347	rBV2	22463	32801	4.11%	0.415%
89	15.408	2347	2349	2353	rVV3	12296	15523	1.95%	0.197%
90	15.481	2356	2361	2369	rVV3	40744	80907	10.14%	1.025%
91	15.566	2370	2375	2379	rVV	60110	99550	12.48%	1.261%
92	15.615	2379	2383	2386	rVV	66075	101542	12.73%	1.286%
93	15.652	2386	2389	2397	rVB2	45042	73332	9.19%	0.929%
94	15.737	2399	2403	2408	rBV7	8161	17754	2.23%	0.225%
95	15.847	2413	2421	2433	rVB2	27063	83259	10.44%	1.054%
96	15.987	2438	2444	2450	rBV3	26119	59907	7.51%	0.759%
97	16.097	2457	2462	2470	rBV2	41834	76172	9.55%	0.965%
98	16.554	2535	2537	2540	rBV3	6728	9058	1.14%	0.115%
99	16.603	2543	2545	2549	rVV3	17995	23316	2.92%	0.295%
100	16.658	2550	2554	2560	rVB2	15426	28192	3.53%	0.357%

Sum of corrected areas: 7895849

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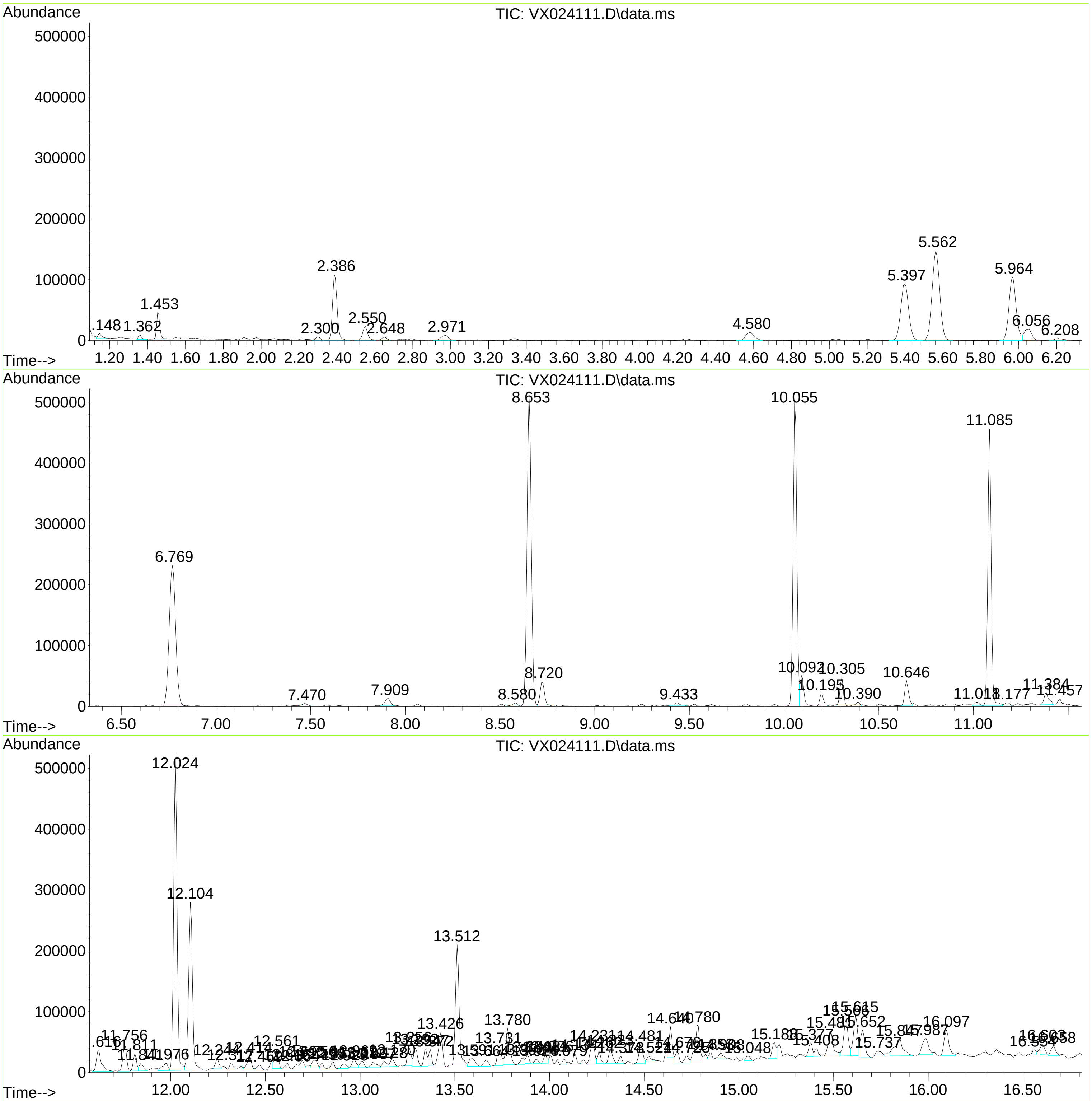
Instrument :
MSVOA_X
ClientSampleId :
1992

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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Data File : VX024111.D
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Instrument :
MSVOA_X
ClientSampleId :
1992

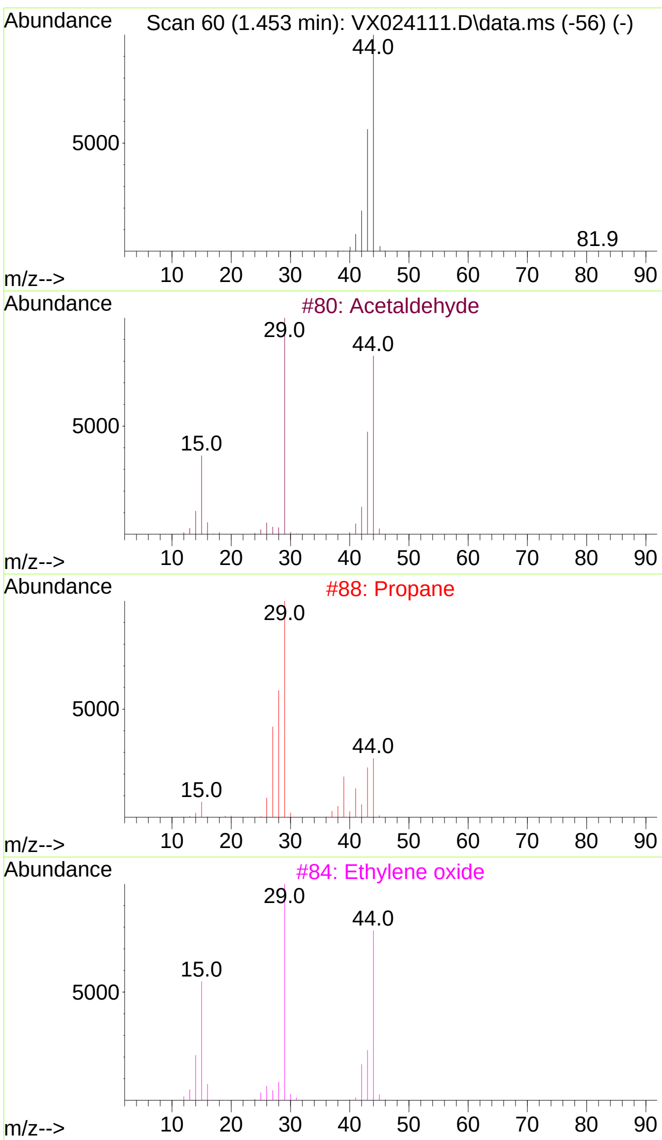
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

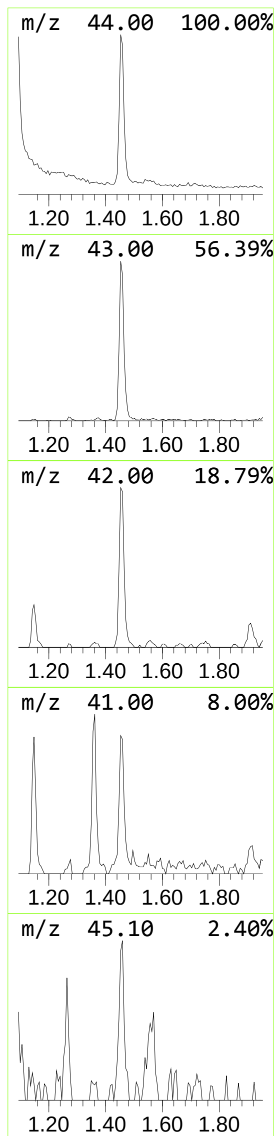
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown1.453 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.453	6.12 ug/l	50623	Pentafluorobenzene	5.562

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetaldehyde	44	C2H4O	000075-07-0	9
2	Propane	44	C3H8	000074-98-6	5
3	Ethylene oxide	44	C2H4O	000075-21-8	5
4	Acetic acid, [(aminocarbonyl)ami...	132	C3H4N2O4	000585-05-7	4
5	Acetic acid, 2-(methylaminoethyl...	117	C5H11NO2	026921-43-7	4





Library Search Compound Report

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Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090321\
Data File : VX024111.D
Acq On    : 03 Sep 2021  18:37
Operator  : JC/MD
Sample    : M3646-13
Misc      : 5.0mL/MSVOA_X/WATER
ALS Vial  : 23   Sample Multiplier: 1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title  : SW846 8260

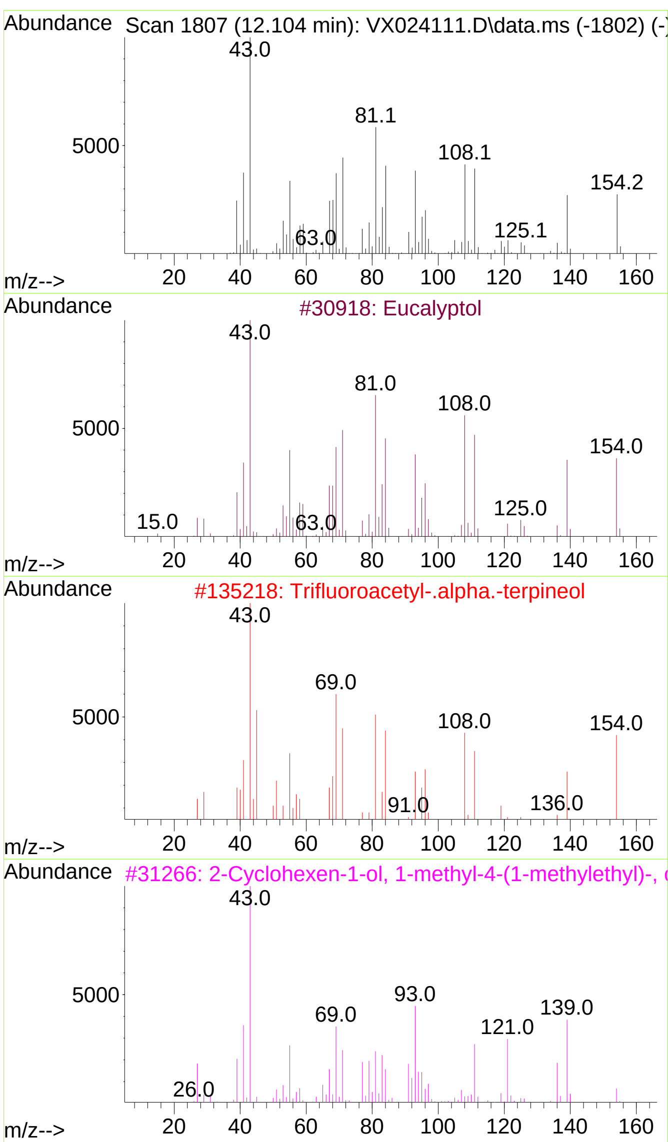
TIC Library   : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P
```

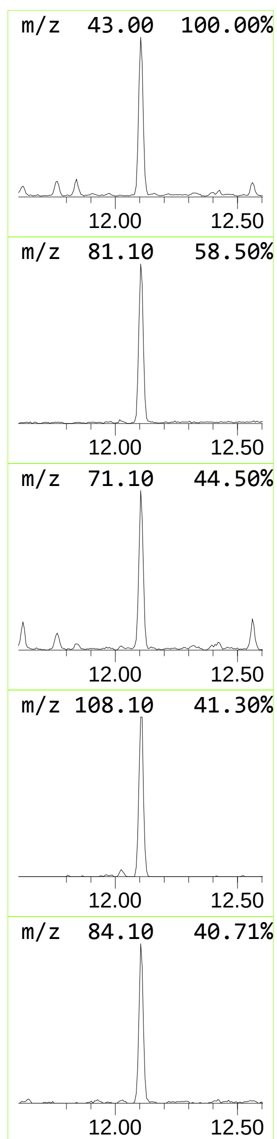
Instrument :
MSVOA_X
ClientSampleId :
1992

```
*********************************************************************
Peak Number  2  Eucalyptol                      Concentration Rank  1
```

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.104	30.95 ug/l	397066	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eucalyptol	154	C10H18O	000470-82-6	99
2	Trifluoroacetyl-.alpha.-terpineol	250	C12H17F3O2	1000058-17-6	35
3	2-Cyclohexen-1-ol, 1-methyl-4-(1...	154	C10H18O	029803-82-5	27
4	2-Isopropyl-5-methylhex-2-enal	154	C10H18O	035158-25-9	22
5	2-Acetylcyclopentanone	126	C7H10O2	001670-46-8	16





Instrument :
MSVOA_X
ClientSampleId :
1992

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Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090321\
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Operator  : JC/MD
Sample    : M3646-13
Misc      : 5.0mL/MSVOA_X/WATER
ALS Vial  : 23   Sample Multiplier: 1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title  : SW846 8260

TIC Library    : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P
```

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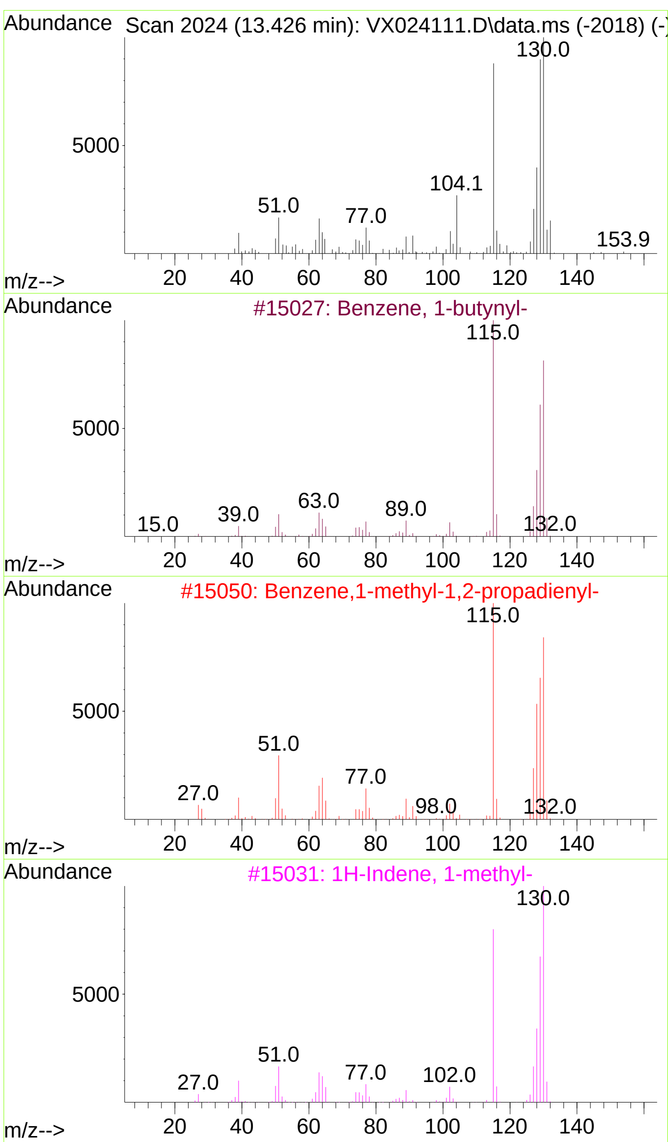
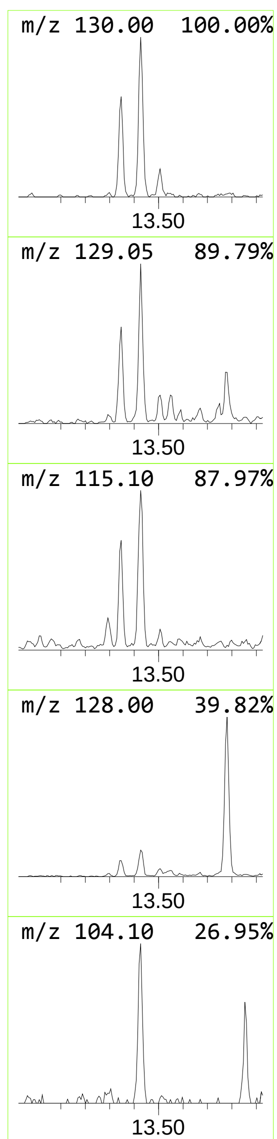
Peak Number 3 Benzene, 1-butynyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.427	6.72 ug/l	86158	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-butynyl-	130	C10H10	000622-76-4	95
2	Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	93
3	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	91
4	2a,4a,6a,6b-Tetrahydrocyclopenta...	130	C10H10	006053-74-3	90
5	Naphthalene, 1,2-dihydro-	130	C10H10	000447-53-0	90





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Instrument :
MSVOA_X
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

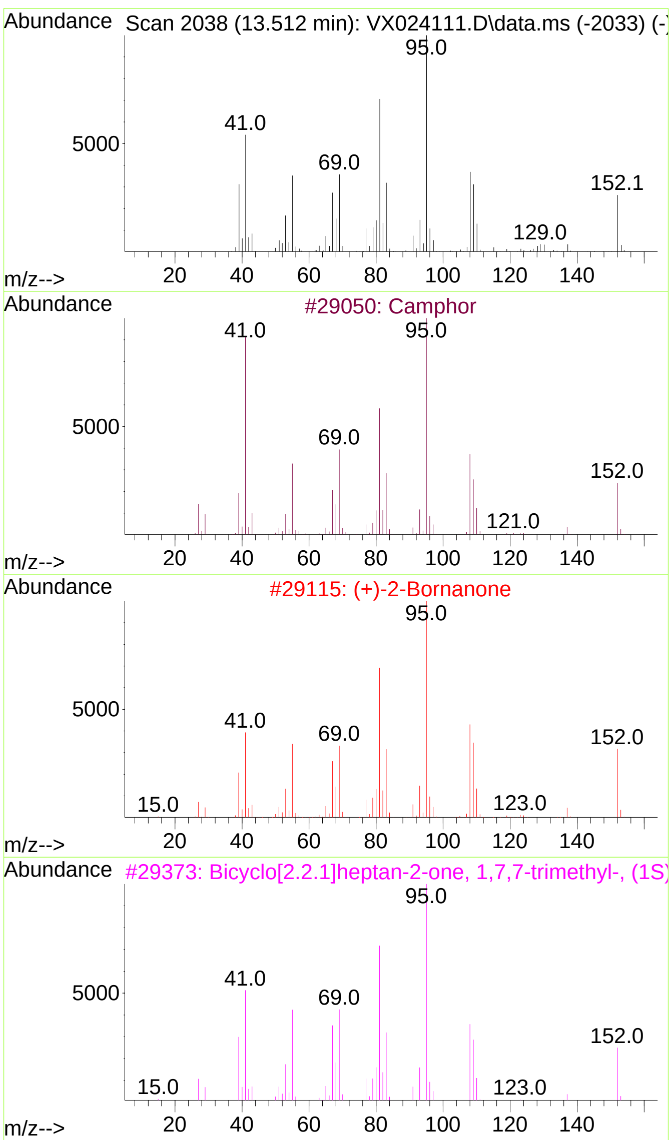
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

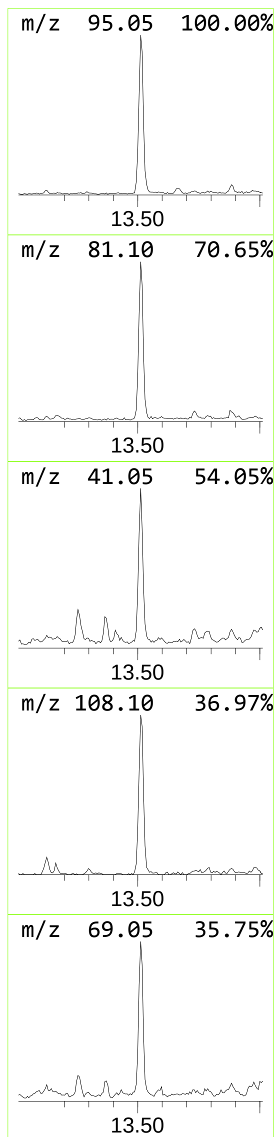
****

Peak Number 4 Camphor Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.512	20.50 ug/l	262971	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Camphor	152	C10H16O	000076-22-2	98
2	(+)-2-Bornanone	152	C10H16O	000464-49-3	98
3	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	98
4	2-Butanone, 4-(5-methyl-2-furanyl)-	152	C9H12O2	013679-56-6	49
5	Tricyclo[2.2.1.0(2,6)]heptan-3-o...	152	C10H16O	062560-53-6	49





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ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
1992

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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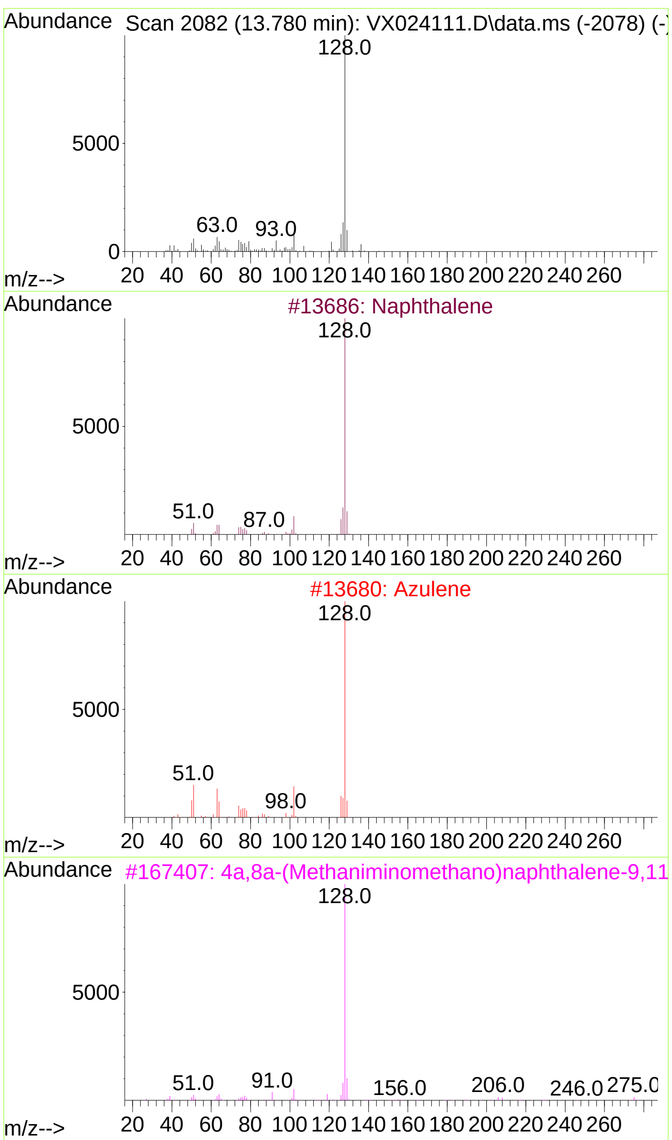
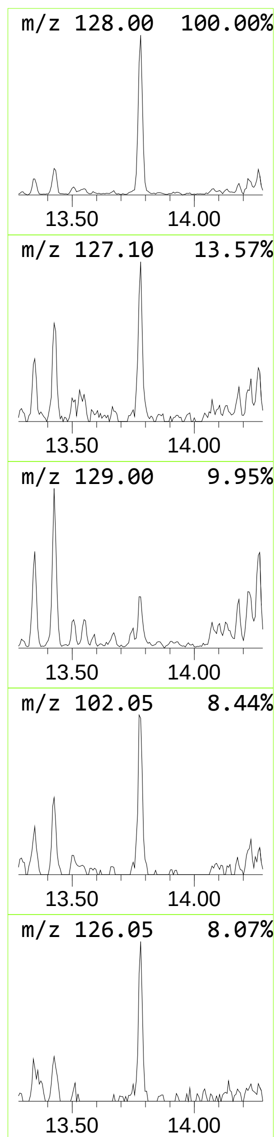
Peak Number 5 Azulene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.780	6.99 ug/l	89606	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	94
2	Azulene	128	C10H8	000275-51-4	90
3	4a,8a-(Methaniminomethano)naphth...	275	C18H13NO2	069915-10-2	72
4	1H-Indene, 1-methylene-	128	C10H8	002471-84-3	47
5	N-Methyl-9-aza-tricyclo[6.2.2.0(...	185	C12H11NO	013131-19-6	45





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ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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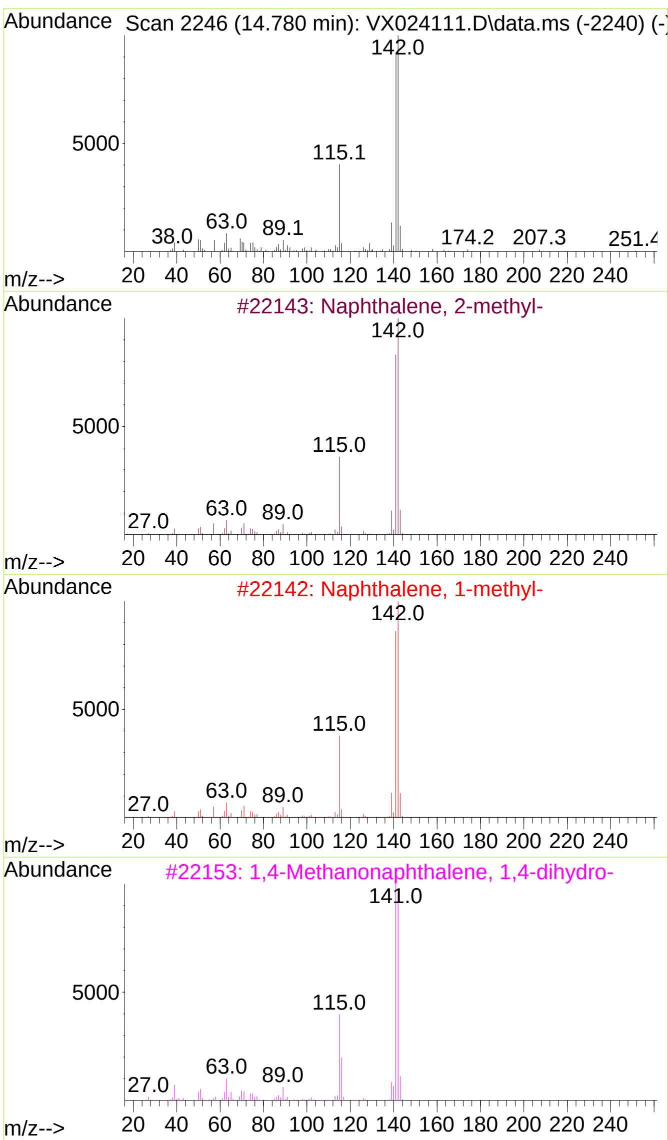
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

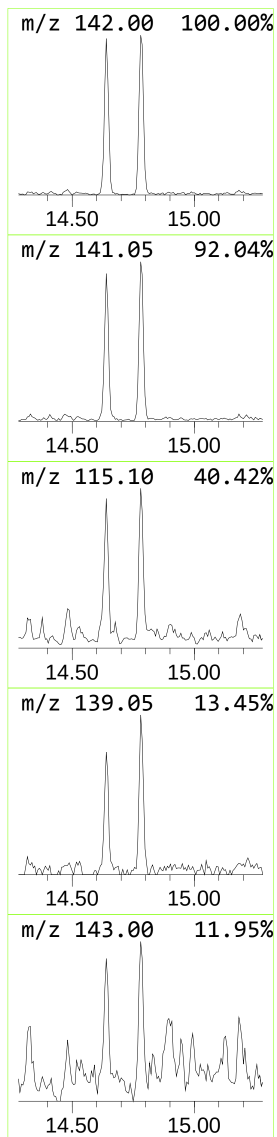
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	7.09 ug/l	90903	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090321\
Data File : VX024111.D
Acq On : 03 Sep 2021 18:37
Operator : JC/MD
Sample : M3646-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
1992

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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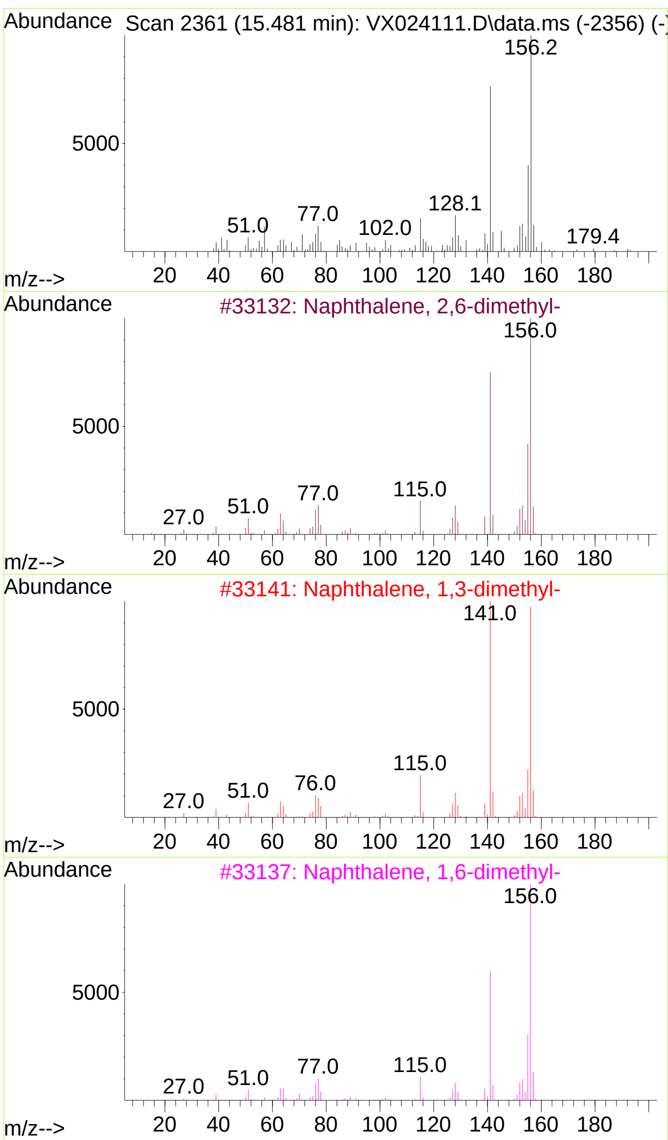
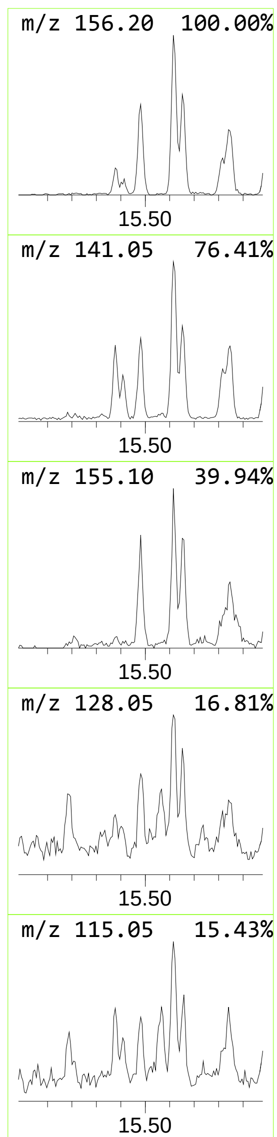
Peak Number 7 Naphthalene, 2,6-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.481	6.31 ug/l	80907	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
2	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
4	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96
5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96





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Acq On : 03 Sep 2021 18:37
Operator : JC/MD
Sample : M3646-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
1992

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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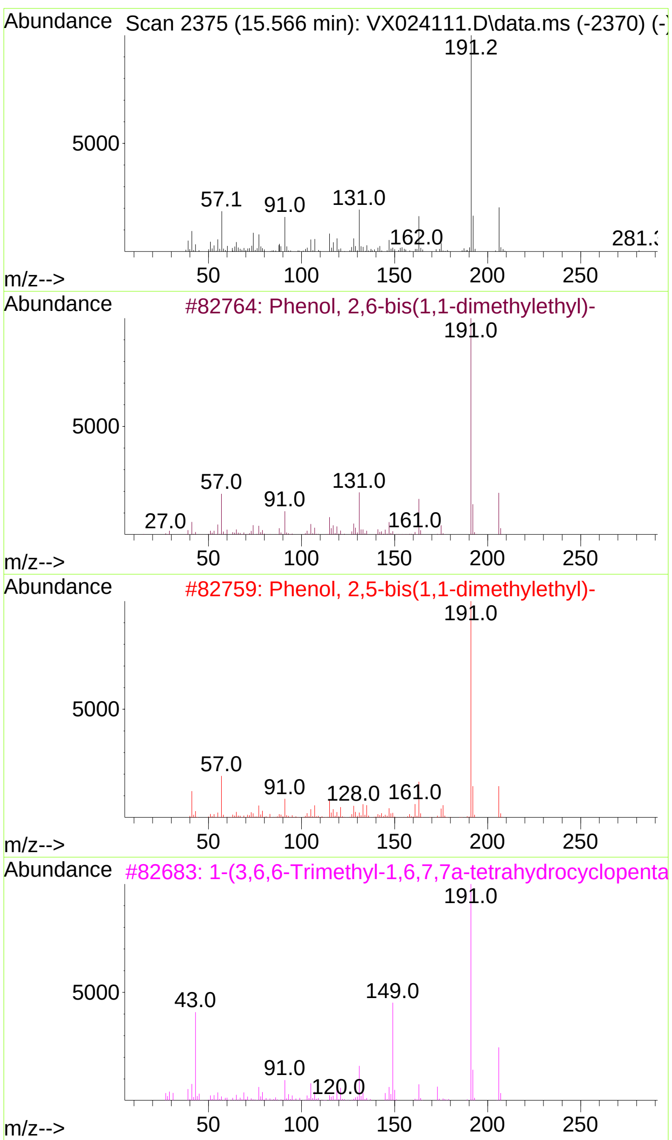
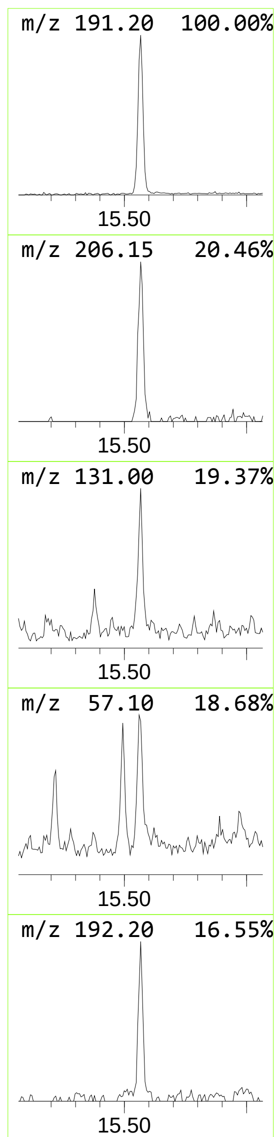
Peak Number 8 Phenol, 2,6-bis(1,1-dimethy... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.566	7.76 ug/l	99550	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-bis(1,1-dimethylethyl)-	206	C14H22O	000128-39-2	96
2	Phenol, 2,5-bis(1,1-dimethylethyl)-	206	C14H22O	005875-45-6	87
3	1-(3,6,6-Trimethyl-1,6,7,7a-tetr...	206	C13H18O2	1000194-97-2	83
4	2,4-Di-tert-butylphenol	206	C14H22O	000096-76-4	72
5	Phenol, 3,5-bis(1,1-dimethylethyl)-	206	C14H22O	001138-52-9	68





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Data File : VX024111.D
Acq On : 03 Sep 2021 18:37
Operator : JC/MD
Sample : M3646-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
1992

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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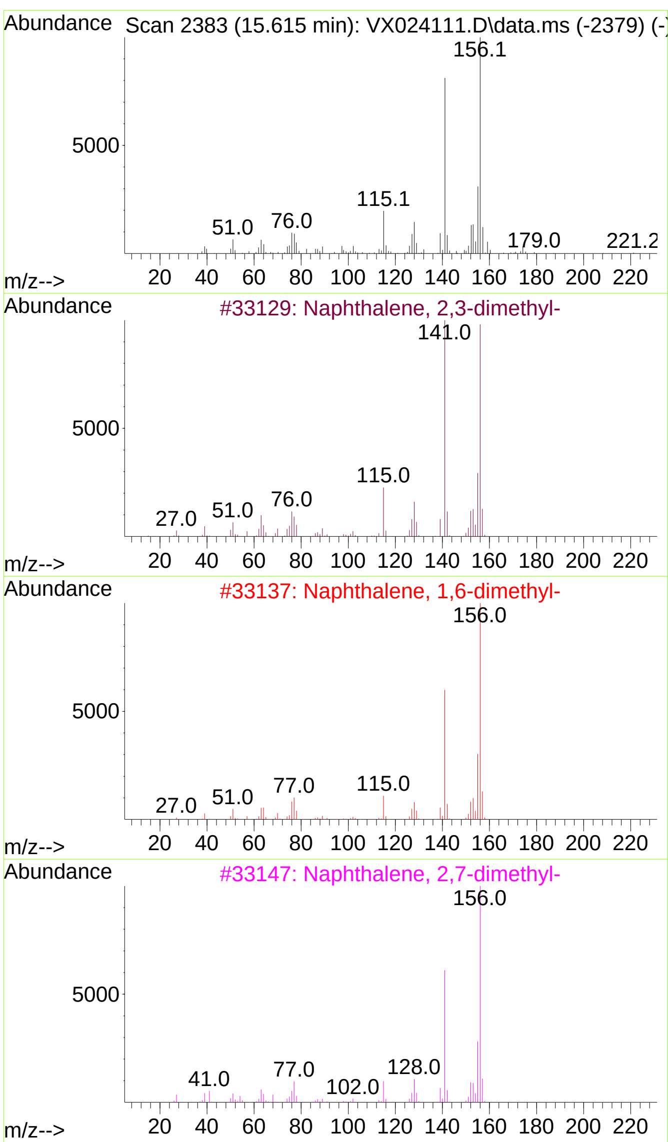
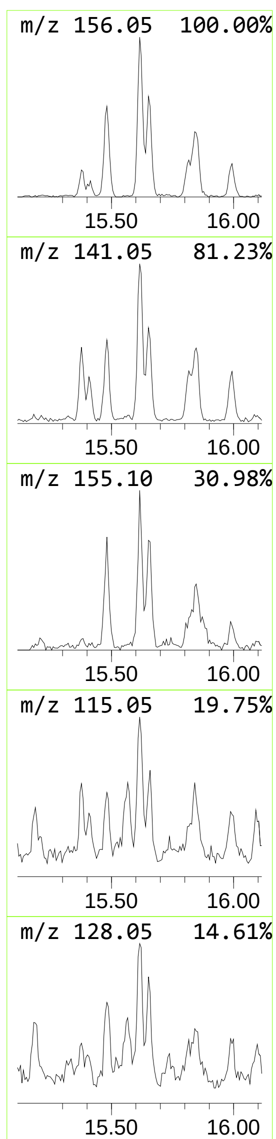
Peak Number 9 Naphthalene, 2,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.615	7.92 ug/l	101542	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96





Library Search Compound Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090321\
Data File : VX024111.D
Acq On : 03 Sep 2021 18:37
Operator : JC/MD
Sample : M3646-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
1992

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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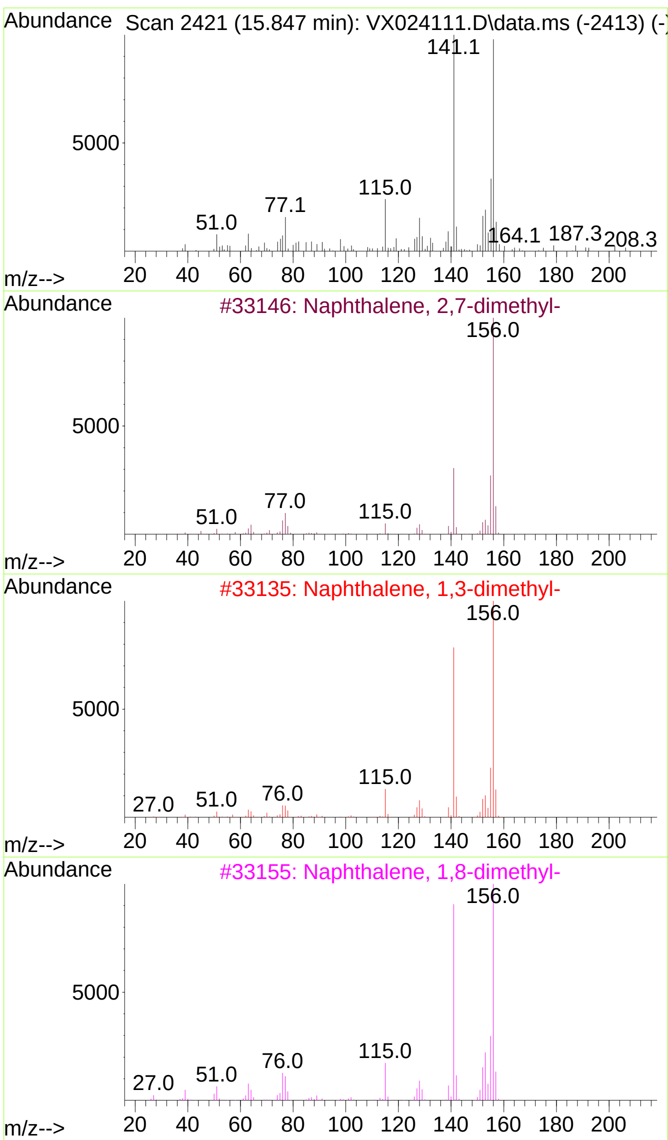
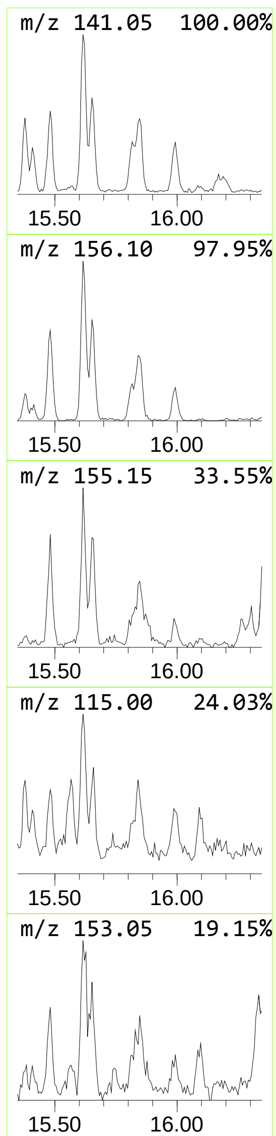
Peak Number 10 Naphthalene, 2,7-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.847	6.49 ug/l	83259	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
2	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	94
3	Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	93
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	93
5	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090321\
Data File : VX024111.D
Acq On : 03 Sep 2021 18:37
Operator : JC/MD
Sample : M3646-13
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
1992

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
unknown1.453	1.453	6.1	ug/l	50623	1	5.562	413858	50.0
Eucalyptol	12.104	30.9	ug/l	397066	4	12.024	641364	50.0
Benzene, 1-buty...	13.427	6.7	ug/l	86158	4	12.024	641364	50.0
Camphor	13.512	20.5	ug/l	262971	4	12.024	641364	50.0
Azulene	13.780	7.0	ug/l	89606	4	12.024	641364	50.0
Naphthalene, 1-...	14.780	7.1	ug/l	90903	4	12.024	641364	50.0
Naphthalene, 2,...	15.481	6.3	ug/l	80907	4	12.024	641364	50.0
Phenol, 2,6-bis...	15.566	7.8	ug/l	99550	4	12.024	641364	50.0
Naphthalene, 2,...	15.615	7.9	ug/l	101542	4	12.024	641364	50.0
Naphthalene, 2,...	15.847	6.5	ug/l	83259	4	12.024	641364	50.0